result in loss of user privileges and other penalties.

FILE 'HOME' ENTERED AT 14:23:12 ON 20 NOV 2006

=> fil reg
COST IN U.S. DOLLARS

SINCE FILE TOTAL
ENTRY SESSION
0.21 0.21

FULL ESTIMATED COST

FILE 'REGISTRY' ENTERED AT 14:23:22 ON 20 NOV 2006
USE IS SUBJECT TO THE TERMS OF YOUR STN CUSTOMER AGREEMENT.
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Property values tagged with IC are from the ZIC/VINITI data file provided by InfoChem.

STRUCTURE FILE UPDATES: 19 NOV 2006 HIGHEST RN 913611-00-4 DICTIONARY FILE UPDATES: 19 NOV 2006 HIGHEST RN 913611-00-4

New CAS Information Use Policies, enter HELP USAGETERMS for details.

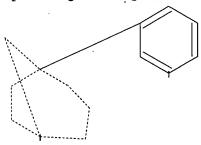
TSCA INFORMATION NOW CURRENT THROUGH June 30, 2006

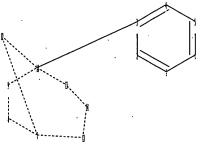
Please note that search-term pricing does apply when conducting SmartSELECT searches.

REGISTRY includes numerically searchable data for experimental and predicted properties as well as tags indicating availability of experimental property data in the original document. For information on property searching in REGISTRY, refer to:

http://www.cas.org/ONLINE/UG/regprops.html

Uploading C:\Program Files\Stnexp\Queries\10-500015z.str





ring nodes :

1 2 3 4 5 6 7 8 9 10 11 12 13 14

chain bonds :

3-10

ring bonds :

1-2 1-6 2-3 3-4 4-5 5-6 7-8 7-12 7-13 8-9 9-10 10-11 10-12 11-14

13-14

exact/norm bonds :

7-8 7-12 7-13 8-9 9-10 10-11 10-12 11-14 13-14

exact bonds :

3-10

normalized bonds :

1-2 1-6 2-3 3-4 4-5 5-6

Match level :

1:Atom 2:Atom 3:Atom 4:Atom 5:Atom 6:Atom 7:Atom 8:Atom 9:Atom 10:Atom 11:Atom 12:Atom 13:Atom 14:Atom

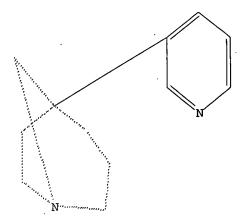
L1 STRUCTURE UPLOADED

=> d

L1 HAS NO ANSWERS

L1

STR



Structure attributes must be viewed using STN Express query preparation.

=> s l1 ful

FULL SEARCH INITIATED 14:23:59 FILE 'REGISTRY'

FULL SCREEN SEARCH COMPLETED - 93 TO ITERATE

100.0% PROCESSED

93 ITERATIONS

70 ANSWERS

SEARCH TIME: 00.00.01

L2

70 SEA SSS FUL L1

=> fil caplus

COST IN U.S. DOLLARS

SINCE FILE

TOTAL

FULL ESTIMATED COST

ENTRY

SESSION

166.94

167.15

FILE 'CAPLUS' ENTERED AT 14:24:04 ON 20 NOV 2006
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FILE COVERS 1907 - 20 Nov 2006 VOL 145 ISS 22 FILE LAST UPDATED: 19 Nov 2006 (20061119/ED)

Effective October 17, 2005, revised CAS Information Use Policies apply. They are available for your review at:

http://www.cas.org/infopolicy.html

=> s 12

L3 1 L2

=> d fbib abs

```
ANSWER 1 OF 1 CAPLUS COPYRIGHT 2006 ACS on STN
AN 2003:534333 CAPLUS
DN 139:101039
TD perivatives of 5-(pyridin-3-y1)-1-azabicyclo[3.2.1]octane, their preparation, and their application in therapy as nicotinic receptor ligands for treatment of CNS disorders
IN Galli, Frederic; Leclerc, Odile; Lochead, Alistair
PA Sanofi-Synthelabo, Fr.
SO Fr. Demande, 20 pp.
CODEM: FRXXBL
DT Patent
LA French
FRAN.CNT 1
PATENT NO. KIND DATE APPLICATION NO. DATE
           FR 2834511
FR 2834511
CA 2471628
                                                                                      20030711
20040213
20030717
                                                                                                                      FR 2002-109
                                                                                                                                                                                   20020107
              MARPAT 139:101039
```

ANSWER 1 OF 1 CAPLUS COPYRIGHT 2006 ACS on STN (Continued)

$$\bigcap_{N} \bigcap_{i=1}^{R}$$

Title compds. I and their acid addition salts are disclosed [wherein: R = halo, a Ph group (substituted by one or more groups chosen from halo,

alkyl or alkoxy, NO2, amino, CF3, cyano, OH, acetyl, or methylenedioxy), pyridinyl, thienyl, indolyl, or pyrimidinyl (possibly substituted by one or more C1-6 alkoxy); dashed bonds = one single bond and another single

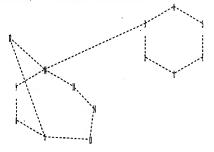
double bond). The compds, are useful as pharmaceuticals, particularly as CNS agents, and specifically as ligands of nicotinic receptors. The compds, were tested against nicotinic receptors with the $\alpha\beta 2$ subunit or with the α^2 subunit. Four synthetic examples and a list of 35 specific compds. (as either di- or tri-HBr or 1:1 oxalate salts)

given. For instance, 2,5-dibromopyridine was arylated in the 2-position by PhB(OH)2 using Pd(PPh3)4 catalyst, and the resultant 5-bromo-2-phenylpyridine was lithiated with Buli and treated with 1-azabicyclo[2.2.2]octan-3-one to give the bicyclic alc. II. Dehydration and rearrangement of II by heating with MeSO3H at 180° gave invention compound III, isolated as the di-HBr salt. In tests for specific binding to isolated rat cerebral nicotinic receptors having either aqR2 or a7 subunits, compds. I had IC50 values in the ranges of 0.01-10 µM and 0.005-20 µM, resp. Some compds. showed selectivity for the a7 receptor subtype.

RE.CNT 3 THERE ARE 3 CITED REFERENCES AVAILABLE FOR THIS RECORD ALL CITATIONS AVAILABLE IN THE RE FORMAT

=>
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ring nodes:
1 2 3 4 5 6 7 8 9 10 11 12 13 14

chain bonds :

3-10

ring bonds :

1-2 1-6 2-3 3-4 4-5 5-6 7-8 7-12 7-13 8-9 9-10 10-11 10-12 11-14

13-14

exact/norm bonds :

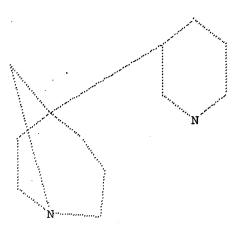
1-2 1-6 2-3 3-4 3-10 4-5 5-6 7-8 7-12 7-13 8-9 9-10 10-11 10-12 11-14 13-14

Match level :

1:Atom 2:Atom 3:Atom 4:Atom 5:Atom 6:Atom 7:Atom 8:Atom 9:Atom 10:Atom 11:Atom 12:Atom 13:Atom 14:Atom

L4 STRUCTURE UPLOADED

=> d L4 HAS NO ANSWERS L4 STR



Structure attributes must be viewed using STN Express query preparation.

=> s 14 ful
 REG1stRY INITIATED
Substance data SEARCH and crossover from CAS REGISTRY in progress...
Use DISPLAY HITSTR (or FHITSTR) to directly view retrieved structures.

FULL SEARCH INITIATED 14:25:34 FILE 'REGISTRY'
FULL SCREEN SEARCH COMPLETED - 403 TO ITERATE

100.0% PROCESSED

403 ITERATIONS

70 ANSWERS

SEARCH TIME: 00.00.01

L5

70 SEA SSS FUL L4

L6

1 L5

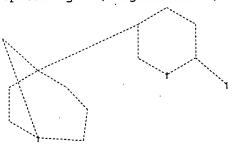
=>

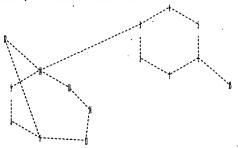
=> d bib

Page 9

=>

Uploading C:\Program Files\Stnexp\Queries\10-500015zb.str





chain nodes :

15

ring nodes :

1 2 3 4 5 6 7 8 9 10 11 12 13 14

chain bonds :
3-10 6-15
ring bonds :

1-2 1-6 2-3 3-4 4-5 5-6 7-8 7-12 7-13 8-9 9-10 10-11 10-12 11-14

13-14

exact/norm bonds :

 $1-2 \quad 1-6 \quad 2-3 \quad 3-4 \quad 3-10 \quad 4-5 \quad 5-6 \quad 6-15 \quad 7-8 \quad 7-12 \quad 7-13 \quad 8-9 \quad 9-10 \quad 10-11 \quad 10-12$

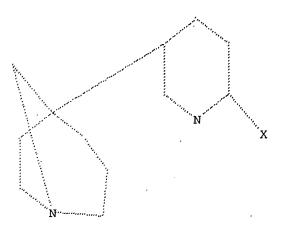
11-14 13-14

Match level :

1:Atom 2:Atom 3:Atom 4:Atom 5:Atom 6:Atom 7:Atom 8:Atom 9:Atom 10:Atom 11:Atom 12:Atom 13:Atom 14:Atom 15:CLASS

L7 STRUCTURE UPLOADED

=> d L7 HAS NO ANSWERS L7 STR



Structure attributes must be viewed using STN Express query preparation.

=> s 17 ful

REG1stRY INITIATED

Substance data SEARCH and crossover from CAS REGISTRY in progress... Use DISPLAY HITSTR (or FHITSTR) to directly view retrieved structures.

FULL SEARCH INITIATED 14:45:16 FILE 'REGISTRY'
FULL SCREEN SEARCH COMPLETED - 9 TO ITERATE

100.0% PROCESSED

9 ITERATIONS

4 ANSWERS

SEARCH TIME: 00.00.01

1.8

4 SEA SSS FUL L7

L9

1 L8

=> d bib